

### Tools for top-down performance analysis of GPUaccelerated applications

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#### **GPU Performance Tools**



- Trace view
  - A series of events that happen over time on each process, thread, and GPU stream
- Profile view
  - A correlation of performance metrics with program contexts
- Existing GPU performance tools
  - GPU vendors
    - Nsight Systems, Nsight Compute, nvprof, ROCProfiler, Intel VTune
  - Third parties
    - TAU, VampirTrace, Allinea Map





- Existing performance tools are ill-suited for analyzing complex programs because they lack a comprehensive profile view to analyze
  - Sophisticated CPU calling contexts
    - A GPU API (e.g., cudaMemcpy) invoked in different places
  - Sophisticated GPU calling contexts
    - OpenMP Target, Kokkos, and RAJA generate code with many small procedures
- At best, existing tools only attribute runtime cost to a flat profile view of functions executed on GPUs



#### **HPCToolkit Profile View**



RuclearData.cc 🕴

246// Return the total cross section for this energy group 247 HOST DEVICE 248 double NuclearData::getReactionCrossSection( unsigned int reactIndex, unsigned int isotopeIndex, unsigned int group) 249 250 { 251 qs\_assert(isotopeIndex < \_isotopes.size());</pre> 252 qs assert(reactIndex < isotopes[isotopeIndex]. species[0]. reactions.size());</pre> 253 return isotopes[isotopeIndex]. species[0]. reactions[reactIndex].getCrossSection(group); 254 }

🕆 Top-down view 🖾 🗞 Bottom-up view 👬 Flat view

🕆 🕂 🍈 fox 🚺 🐖 🗛 👘 🖬 👻

Scope		GINS:Sum (I)	GINS:STL_ANY:Sum (I)
Joop at main.cc: 159		1.07e+11 100 %	9.83e+10 100 %
loop at main.cc: 163		1.07e+11 100 %	9.83e+10 100 %
✓ ➡ 193: [I] CycleTrackingKernel(MonteCarlo*, int, ParticleVault*, ParticleVault*)		1.07e+11 100 %	9.83e+10 100 %
🧼 🖶 127:device_stubZ19CycleTrackingKernelP10MonteCarloiP13ParticleVaultS2_(MonteCarlo*, int, Parti		1.07e+11 100 😵	9.83e+10 100 <mark>%</mark>
CPU Calling Context 🧹 ា 14: [I] cudaLaunchKernel <char></char>		1.07e+11 100 %	9.83e+10 100 %
GPU API Node 🛛 🗸 🖶	209: <gpu kernel=""></gpu>	1.07e+11 100 %	9.83e+10 100 %
~	174: CycleTrackingKernel(MonteCarlo*, int, ParticleVault*, ParticleVault*)	1.07e+11 100 %	9.83e+10 100 %
B) 132: CycleTrackingGuts(MonteCarlo*, int, ParticleVault*, ParticleVault*)		1.06e+11 100.0	9.82e+10 100.0
loop at CycleTracking.cc: 118		8.90e+10 83.5%	8.08e+10 82.2%
		4.99e+10 46.9%	4.49e+10 45.7%
[I] inlined from QS_Vector.hh: 94 loop at QS_Vector.hh: 94		3.76e+10 35.3%	3.34e+10 34.0%
		3.61e+10 33.9%	3.20e+10 32.5%
	<ul> <li>[I] inlined from CollisionEvent.cc: 71</li> </ul>	3.58e+10 33.6%	3.17e+10 32.3%
	loop at CollisionEvent.cc: 71	3.42e+10 32.1%	3.03e+10 30.9%
GPU Loops and Inline Func	Functions 🔰 🤍 🖶 73: macroscopicCrossSection(MonteCarlo*, int, int, int, int, int)	3.11e+10 29.2%	2.78e+10 28.3%
	<ul> <li>[I] inlined from MacroscopicCrossSection.cc: 45</li> </ul>	2.57e+10 24.1%	2.31e+10 23.5%
	↓ ➡ 41: NuclearData::getReactionCrossSection(unsigned int, unsigned	1.69e+10 15.9%	1.56e+10 15.9%
GPU Calling Context	🧹 🥪 [l] inlined from NuclearData.cc: 194	9.36e+09 8.8%	8.70e+09 8.9%
GPU Hotspot	NuclearData.cc: 253	9.06e+09 8.5%	8.44e+09 P.60
8/21/2020			4

#### Outline



#### Overview

- Performance Measurement
- Performance Metrics Attribution
- Performance Analysis
- Case Studies
- Summary



#### **HPCToolkit Overview**





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#### **HPCToolkit Workflow**









- GPU measurement collection
  - Multiple application threads launching kernels to a GPU
  - A monitor thread reads measurements and attributes them to the corresponding application threads
- GPU API correlation in CPU calling context tree
  - Thousands of GPU invocations, including kernel launches, memory copies, and synchronizations in large-scale applications
- GPU metrics attribution
  - Read line map and DWARF in heterogenous binaries
  - Identify loop nests
  - Reconstruct GPU calling context



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- Two categories of threads
  - Application Threads (*N* per process)
    - Launch kernels, move data, and synchronize GPU calls
  - Monitor Thread (1 per process)
    - Monitor GPU events and collect GPU measurements
- Interaction
  - **Create correlation:** An application thread *T* creates a correlation record when it launches a kernel and tags the kernel with a correlation ID *C*, notifying the monitor thread that *C* belongs to *T*
  - Attribute measurements: The monitor thread collects measurements associated with *C* and communicates measurement records back to thread *T*





- Communication channels: wait-free unordered stack groups
- A private stack and a shared stack used by two threads
  - **Producer PUSH**(*CAS*): push an item on a shared stack
  - **Consumer STEAL**(*XCHG*): steal the contents of the shared stack, push the contents onto a private stack
  - **Consumer POP**: pop an item from the private stack
- Wait-free because PUSH fails at most once when a concurrent thread STEALs contents of the shared stack



### Interactions between the GPU monitor thread and application threads







# GPU API Correlation with CPU Calling Context



- Unwind a call stack from each API invocation, including kernel launch, memory copy, and synchronization
- Initial approach:
  - Identify the function enclosing each call site in the call stack using a global shared map
- Problem:
  - Applications have deep call stacks and large codebase
    - Nyx: up to 60 layers and 400k calls
    - Laghos: up to 40 layers and 100k calls



#### Fast Unwinding



- Refined approach:
  - Memoize common call path prefixes
    - Temporally-adjacent samples in complex applications often share common call path prefixes
    - Employ eager (mark bits) or lazy (tramopoline) marking to identify LCA of call stack unwinds
  - Avoid costly access to mutable concurrent data
    - Cache unwinding recipes in a per thread hash table
  - Avoid duplicate unwinds
    - Filter CUDA Driver APIs within CUDA Runtime APIs



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- Attribute metrics to flat PCs at runtime
  - Relocate each GPU function in a CUBIN so that they do not overlap
- Aggregate metrics to lines
  - Read .debug\_lineinfo section if available
- Aggregate metrics to loops
  - nvdisasm –poff –cfg for all valid functions
  - Parse dot files to data structures for Dyninst
  - Use Dyninst ParseAPI to identify loops



#### **GPU Calling Context Tree**



- Problem
  - Unwinding call stacks on GPU is costly for each GPU thread
  - NVIDIA's CUPTI does not provide an unwinding API
- Challenges
  - GPU functions may be invoked from different call sites
  - Need to decide how to attribute costs to each call site
- Solution
  - Reconstruct GPU calling context tree from flat instruction samples and static GPU call graph



#### Reconstruct Approximate GPU Calling Context Tree



- Construct static call graph
  - Link call instructions with corresponding functions
- Construct dynamic call graph
  - Propagate call instructions to all possible call sites
  - Prune functions with no samples or calls
- Transform call graph to calling context tree
  - Apportion each function's samples based on samples of its incoming call sites
- See the paper for how we handle recursive calls



GPU Calling Context Tree Example



 D is split into D' and D' based on (0x30, 1) and (0x40, 2)





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#### **CPU** Importance



• CPU<sub>IMPORTANCE</sub>:

Ratio of a procedure's time to the whole execution time

$$\operatorname{Max}\left(\frac{\operatorname{CPU}_{\text{TIME}} - \operatorname{SUM}(\operatorname{GPU}_{\text{API}_{\text{TIME}}})}{\operatorname{CPU}_{\text{TIME}}}, 0\right) \times \frac{\operatorname{CPU}_{\text{TIME}}}{\operatorname{EXECUTION}_{\text{TIME}}}$$

Ratio of a procedure's pure CPU time. If more time is spent on GPU than CPU, the ratio is set to 0

#### • GPU\_API<sub>TIME</sub>:

- KERNEL<sub>TIME</sub>: cudaLaunchKernel, cuLaunchKernel
- MEMCPY<sub>TIME</sub>: *cudaMemcpy*, *cudaMemcpyAsync*
- MEMSET<sub>TIME</sub>: cudaMemset
- ...



#### **GPU API Importance**



• GPU\_API<sub>IMPORTANCE</sub>

GPU\_API<sub>TIME</sub> SUM(GPU\_API<sub>TIME</sub>)

Consider the importance of the memory copy to all the GPU time

- Find which type of GPU API is the most expensive
  - Kernel: optimize specific kernels with PC Sampling profiling
  - Other APIs: apply optimizations based on calling context





- Map opcodes and modifiers to instruction classes
- Memory ops
  - class.[memory hierarchy].width
- Compute ops
  - class.[precision].[tensor].width
- Control ops
  - class.control.type

• . . .





- Problem
  - GPU PC sampling cannot be used in the same pass with metric collection
  - Nsight-compute runs nine passes to collect multiple metrics for a small kernel
- Our approach
  - Derive multiple metrics using PC sampling and other activity records
    - e.g., instruction throughput, scheduler issue rate, SM active ratio
- See the paper for how we derive metrics



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#### **Case Studies**

- Setup
  - Summit compute node: Power9+Volta V100
  - module load hpctoolkit/2020.03.01
  - module load cuda/10.1.243
- Case studies
  - RAJAPerf Suite
    - Performance benchmark for a template-based GPU programming model
  - Laghos
    - A DOE mini-app that solves the time-dependent Euler equation of compressible gas dynamics
  - Nekbone
    - A subset of kernels from Nek5000, a high-order Navier-Stokes solver based on the spectral element method







- Assess the accuracy of our GPU call graph reconstruction
- Compare our call count approximation with exact call counts from NVIDIA's NVBit
  - Compiled with "-G" to avoid function inlining

Test Case	Unique Call Paths	Error
Basic_INIT_VIEW1D_OFFSET	9	0
Basic_REDUCE3_INT	113	0.03
Stream_DOT	60	0.006
Stream_TRIAD	5	0
Apps_PRESSURE	6	0
Apps_FIR	5	0
Apps_DEL_DOT_VEC_2D	3	0
Apps_VOL3D	4	0







- Pinpoint performance problems in profile view by *importance metrics* 
  - CPU takes 80% execution time
    - *mfem::LinearForm::Assemble* only has CPU code, taking 60% execution time
  - Memory copies can be optimized by different methods based on their calling context
    - Use memory copy counts and bytes to determine if using pinned memory with help
    - Eliminate conditional memory copies
    - Fuse memory copies into kernel code



#### Laghos-CUDA



- Original time: 32.9s
  - 11.3s on GPU computation and memory copies
- Optimized time: 30.9s
  - 9.0s on GPU computation and memory copies
- Overall improvement: 6.4%
- GPU code section improvement: 25.6%

#### Laghos-RAJA



- Pinpoint synchronization
  - Kernel launch in CUDA is asynchronous, but Laghos uses RAJA synchronous kernel launch
  - Use asynchronous RAJA kernel launch
- Bad compiler generated code with RAJA template wrapper
  - rMassMultAdd<3,4>: RAJA version has 4x STG instructions as the CUDA version. ¼ STG instructions within a loop use the same address.
  - Store temporary values in local variables



#### Laghos-RAJA



- Original time: 41.0s
  - 19.47s on GPU computation and memory copies
- Optmizied time: 32.2s
  - 10.8s on GPU computation and memory copies
- Overall improvement: 27.3%
- GPU code section improvement: 80.2%







- Associate PC samples with GPU calling context, loops, and lines
- Use instruction mix to provide essential metrics for generating a roofline model
- Problems and optimizations
  - **Memory throttling**: high frequency global memory requests do not always hit cache. +*shared memory*
  - **Memory dependency**: compiler (-O3) does not reorder global memory read properly to hide latency. +*reorder* global memory read
  - Execution dependency: complicated assembly code for integer division. +precompute reciprocal to simplify division





- Overall improvement: +34%
- Estimate errors: the first one +8% because of GPU instruction predicates





- 83% of peak performance
  - Could obtain +19% by fusing multiply and add on the assembly level



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- HPCToolkit pinpoints performance problems for both large-scale applications and individual kernels
- HPCToolkit provides insights for finding problems in compiler-generated GPU code, resource usage, synchronization, parallelism level, instruction pipeline, and memory access patterns
- HPCToolkit collects measurement data efficiently
  - CUDA/10.1
  - Without PC sampling: comparable with nvprof
  - With PC sampling: 6x speedup









### Memoizing common call path prefixes





Eager LCA Arnold & Sweeny, IBM TR, 1999.

Lazy LCA Froyd/@tal, ICS05.

- mark frame RAs while unwinding
- return from marked frame clears mark
- new calls create
   unmarked frame RAs
- mark frame RA during next unwind
- prior marked frames are common prefix

- return from marked frame moves mark
- new calls create unmarked frames
- mark frame RA during next unwind
- prior marked frame indicates common prefix 38



 Link call instructions with corresponding functions





- Challenge
  - Call instructions are sampled (Unlike gprof)
- Assumptions
  - If a function is sampled, it must be called somewhere
  - If there are no call instruction samples for a sampled function, we assign each potential call site one call sample



- Assign call instruction samples to call sites
- Mark a function with *T* if it has instruction samples, otherwise *F*





- Propagate call instructions
  - At the same time change function marks
  - Implemented with a queue





- Prune functions with no samples or calls
- Keep call instructions





- Identify SCCs in call graph
- Link external calls to SCCs and unlink calls inside SCCs



# Step 4: Transform Call Graph to Calling Context Tree



 Apportion each function's samples based on samples of its incoming call sites





- *Stall reason*: When an instruction is sampled, its *stall reason* (if any) is recorded
- Latency sample: If all warps on a scheduler are stalled when a sample is taken, the sample is marked as a *latency sample*



#### Analysis with PC Sampling



- Each stream multiprocessor is sampled individually
- Active warps are uniformly distributed to warp schedulers
- Samples are taken in a fixed number of cycles
- Volta V100
  - scheduler\_id = warp\_index % 4
  - *16* warp slots on each scheduler



Stream Multiprocessor



- Total Samples (S): 6
- Latency Samples (S<sub>L</sub>): 4
- Issue Rate (*I*):  $\frac{S-S_L}{S} = \frac{2}{3}$
- **IPC**: *I* × *MIN*(*active\_warps,warp\_schedulers*)



#### **Estimate Error**



- GPUs are not always running at the maximum clock rate
- Actual average clock rate:  $\bar{C}$
- Maximum clock rate: C
- Estimate error:  $\varepsilon = \frac{c}{\bar{c}}$